

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1623paz

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 Jan 25 BLAST(R) searching in REGISTRY available in STN on the Web
NEWS 3 Jan 29 FSTA has been reloaded and moves to weekly updates
NEWS 4 Feb 01 DKILIT now produced by FIZ Karlsruhe and has a new update
frequency
NEWS 5 Feb 19 Access via Tymnet and SprintNet Eliminated Effective 3/31/02
NEWS 6 Mar 08 Gene Names now available in BIOSIS
NEWS 7 Mar 22 TOXLIT no longer available
NEWS 8 Mar 22 TRCTHERMO no longer available
NEWS 9 Mar 28 US Provisional Priorities searched with P in CA/CAPLUS
and USPATFULL
NEWS 10 Mar 28 LIPINSKI/CALC added for property searching in REGISTRY
NEWS 11 Apr 02 PAPERCHEM no longer available on STN. Use PAPERCHEM2
instead.
NEWS 12 Apr 08 "Ask CAS" for self-help around the clock
NEWS 13 Apr 09 BEILSTEIN: Reload and Implementation of a New Subject Area
NEWS 14 Apr 09 ZDB will be removed from STN
NEWS 15 Apr 19 US Patent Applications available in IFICDB, IFIPAT, and
IFIUDB
NEWS 16 Apr 22 Records from IP.com available in CAPLUS, HCAPLUS, and
ZCAPLUS
NEWS 17 Apr 22 BIOSIS Gene Names now available in TOXCENTER
NEWS 18 Apr 22 Federal Research in Progress (FEDRIP) now available

NEWS EXPRESS February 1 CURRENT WINDOWS VERSION IS V6.0d,
CURRENT MACINTOSH VERSION IS V6.0a(ENG) AND V6.0Ja(JP),
AND CURRENT DISCOVER FILE IS DATED 05 FEBRUARY 2002
NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS INTER General Internet Information
NEWS LOGIN Welcome Banner and News Items
NEWS PHONE Direct Dial and Telecommunication Network Access to STN
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that
specific topic.

All use of STN is subject to the provisions of the STN Customer
agreement. Please note that this agreement limits use to scientific
research. Use for software development or design or implementation
of commercial gateways or other similar uses is prohibited and may
result in loss of user privileges and other penalties.

* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 14:33:48 ON 24 APR 2002

=> file reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.21

0.21

FILE 'REGISTRY' ENTERED AT 14:34:12 ON 24 APR 2002

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2002 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 23 APR 2002 HIGHEST RN 406909-40-8

DICTIONARY FILE UPDATES: 23 APR 2002 HIGHEST RN 406909-40-8

TSCA INFORMATION NOW CURRENT THROUGH July 7, 2001

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES
for more information. See STNote 27, Searching Properties in the CAS
Registry File, for complete details:

<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=>

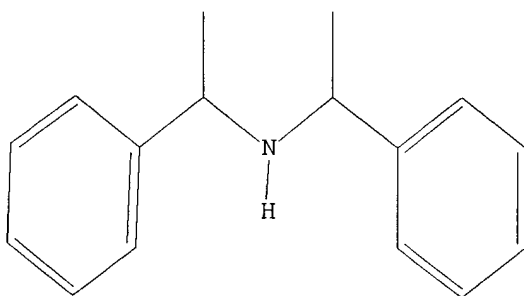
Uploading 09853085 generic amine.str

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> search l1 sss sam

SAMPLE SEARCH INITIATED 14:35:21 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 6695 TO ITERATE

14.9% PROCESSED 1000 ITERATIONS

6 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 129000 TO 138800
PROJECTED ANSWERS: 423 TO 1183

L2 6 SEA SSS SAM L1

=>

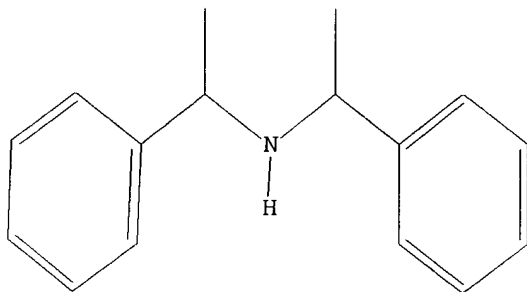
Uploading 09853085 generic amine.str

L3 STRUCTURE UPLOADED

=> d l3

L3 HAS NO ANSWERS

L3 STR



Structure attributes must be viewed using STN Express query preparation.

=> search l3 sss sam

SAMPLE SEARCH INITIATED 14:36:33 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 6695 TO ITERATE

14.9% PROCESSED 1000 ITERATIONS

1 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

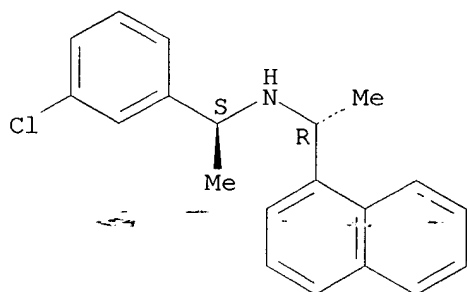
FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 129000 TO 138800
PROJECTED ANSWERS: 1 TO 288

L4 1 SEA SSS SAM L3

=> d scan

L4 1 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 1-Naphthalenemethanamine,
N-[(1S)-1-(3-chlorophenyl)ethyl]-.alpha.-methyl-
, (.alpha.R)- (9CI)
MF C20 H20 Cl N

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> search l3 sss full
 FULL SEARCH INITIATED 14:37:13 FILE 'REGISTRY'
 FULL SCREEN SEARCH COMPLETED - 133082 TO ITERATE

100.0% PROCESSED 133082 ITERATIONS
 SEARCH TIME: 00.00.04

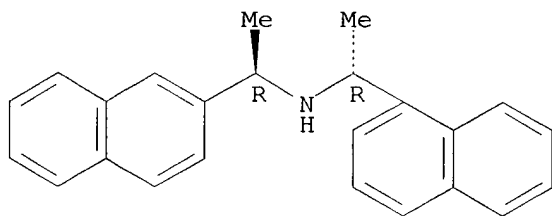
252 ANSWERS

L5 252 SEA SSS FUL L3

=> d scan

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 1-Naphthalenemethanamine,
 .alpha.-methyl-N-[(1R)-1-(2-naphthalenyl)ethyl]-
 , hydrochloride, (.alpha.R)- (9CI)
 MF C24 H23 N . Cl H

Absolute stereochemistry.



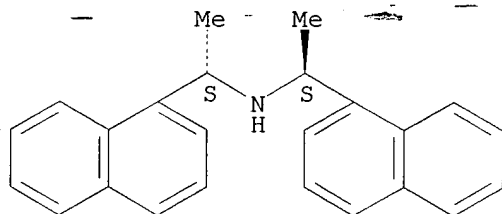
● HCl

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):30

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 1-Naphthalenemethanamine,
 .alpha.-methyl-N-[(1S)-1-(1-naphthalenyl)ethyl]-

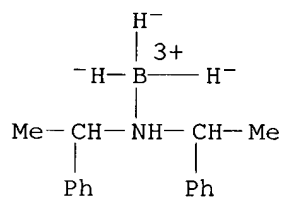
, hydrochloride, (.alpha.S)- (9CI)
 MF C24 H23 N . Cl H

Absolute stereochemistry.



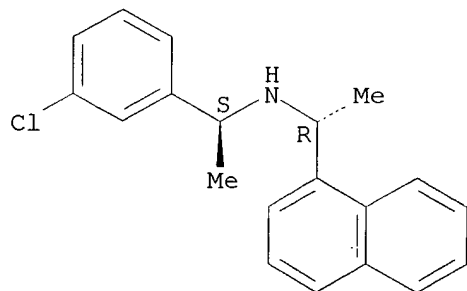
● HCl

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Boron, trihydro[.alpha.-methyl-N-(1-phenylethyl)benzenemethanamine]-,
 [T-4-[S-(R*,R*)]]- (9CI)
 MF C16 H22 B N
 CI CCS



L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 1-Naphthalenemethanamine,
 N-[(1S)-1-(3-chlorophenyl)ethyl]-.alpha.-methyl-
 , (.alpha.R)- (9CI)
 MF C20 H20 Cl N

Absolute stereochemistry.

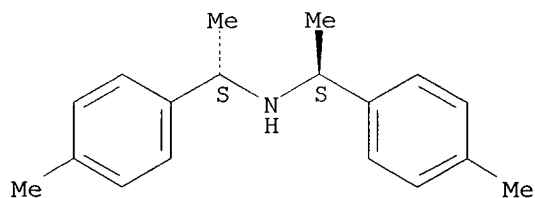


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

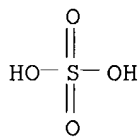
L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenemethanamine, .alpha.,4-dimethyl-N-[(1S)-1-(4-methylphenyl)ethyl]-,
(.alpha.S)-, sulfate (2:1) (9CI)
MF C18 H23 N . 1/2 H2 O4 S

CM 1

Absolute stereochemistry.

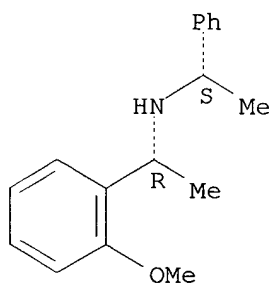


CM 2



L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenemethanamine, 2-methoxy-.alpha.-methyl-N-(1-phenylethyl)-,
[S-(R*,S*)]- (9CI)
MF C17 H21 N O

Absolute stereochemistry.

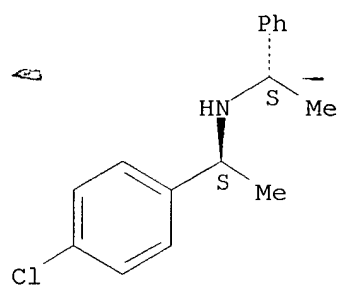


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenemethanamine, 4-chloro-.alpha.-methyl-N-(1-phenylethyl)-, (R*,R*)-

(9CI)
MF C16 H18 Cl N

Relative stereochemistry.

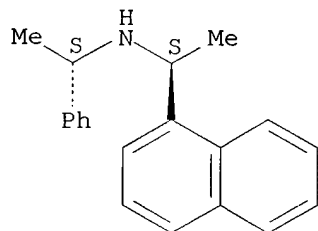


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

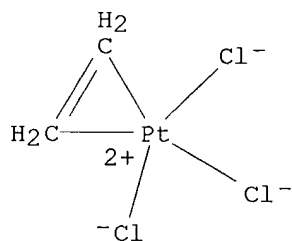
L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Platinate(1-), trichloro(.eta.2-ethene)-, hydrogen, compd. with
(.alpha.S)-.alpha.-methyl-N-[(1S)-1-phenylethyl]-1-naphthalenemethanamine
(1:1) (9CI)
MF C20 H21 N . C2 H4 Cl3 Pt . H

CM 1

Absolute stereochemistry.



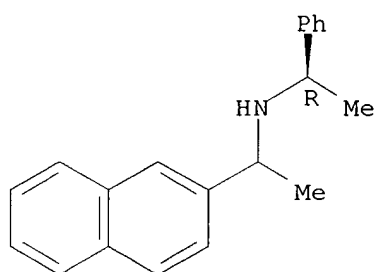
CM 2



H⁺

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2-Naphthalenemethanamine, .alpha.-methyl-N-[(1R)-1-phenylethyl]- (9CI)
 MF C20 H21 N

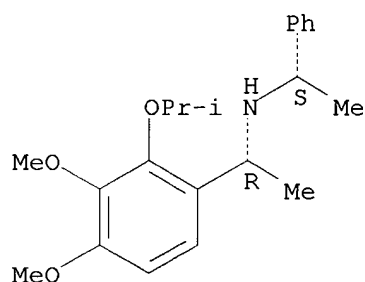
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzenemethanamine, 3,4-dimethoxy-.alpha.-methyl-2-(1-methylethoxy)-N-(1-phenylethyl)-, [S-(R*,S*)]- (9CI)
 MF C21 H29 N O3

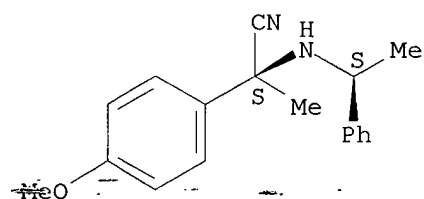
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

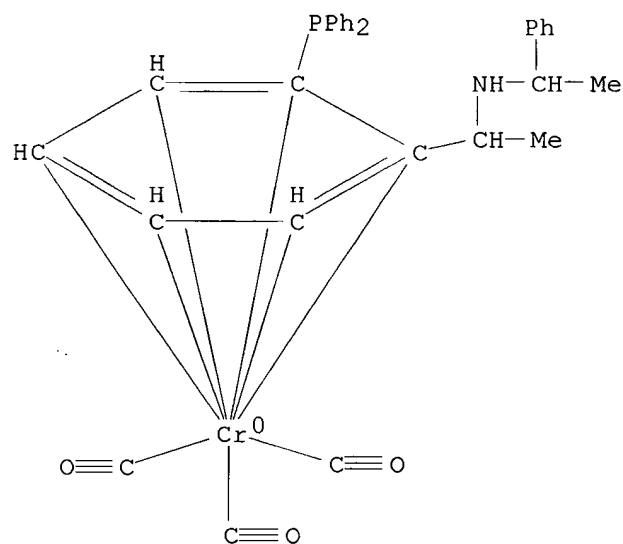
L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzeneacetonitrile, 4-methoxy-.alpha.-methyl-.alpha.-[(1-phenylethyl)amino]-, [S-(R*,R*)]- (9CI)
 MF C18 H20 N2 O

Absolute stereochemistry.



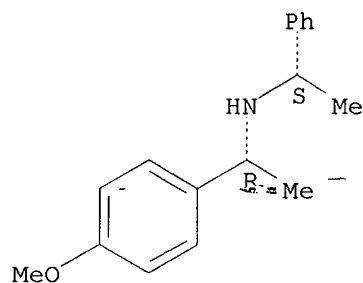
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Chromium, tricarbonyl[(1,2,3,4,5,6-eta.)-2-(diphenylphosphino)-.alpha.-methyl-N-[(1R)-1-phenylethyl]benzenemethanamine]-, stereoisomer (9CI)
 MF C31 H28 Cr N O3 P
 CI CCS



L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzenemethanamine, 4-methoxy-.alpha.-methyl-N-(1-phenylethyl)-, [S-(R*,S*)]- (9CI)
 MF C17 H21 N O

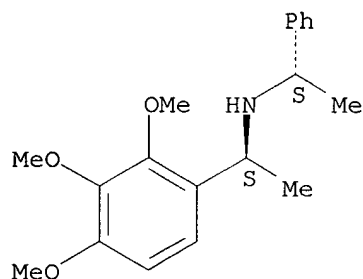
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzenemethanamine, 2,3,4-trimethoxy-.alpha.-methyl-N-(1-phenylethyl)-,
 hydrobromide, [S-(R*,R*)]- (9CI)
 MF C19 H25 N O3 . Br H

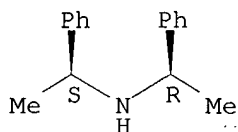
Absolute stereochemistry.



● HBr

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzenemethanamine, .alpha.-methyl-N-[(1R)-1-phenylethyl]-,
 (.alpha.S)-rel- (9CI)
 MF C16 H19 N
 CI COM

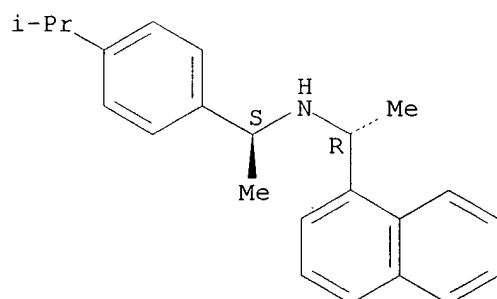
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 1-Naphthalenemethanamine, .alpha.-methyl-N-[(1S)-1-[4-(1-methylethyl)phenyl]ethyl]-, (.alpha.R)- (9CI)
MF C23 H27 N

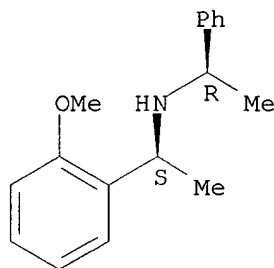
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenemethanamine, 2-methoxy-.alpha.-methyl-N-(1-phenylethyl)-, [R-(R*,S*)]- (9CI)
MF C17 H21 N O

Absolute stereochemistry.

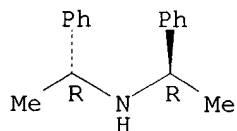


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

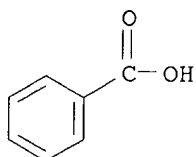
L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenemethanamine, .alpha.-methyl-N-(1-phenylethyl)-, [R-(R*,R*)]-, benzoate (9CI)
MF C16 H19 N . C7 H6 O2

CM 1

Absolute stereochemistry. Rotation (+).

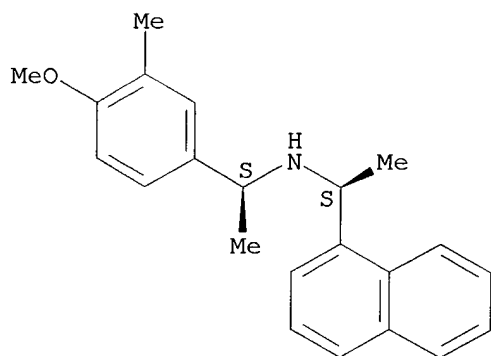


CM 2⁺



L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 1-Naphthalenemethanamine, N-[(1S)-1-(4-methoxy-3-methylphenyl)ethyl]-
 .alpha.-methyl-, (.alpha.S)- (9CI)
 MF C22 H25 N O
 CI COM

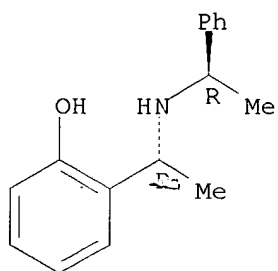
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

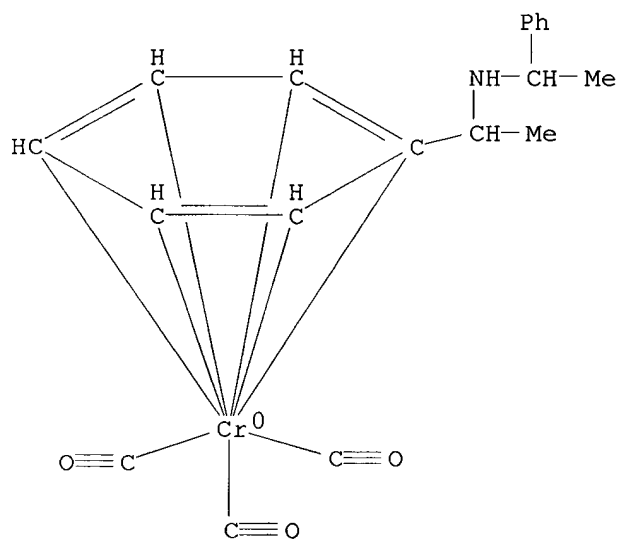
L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Phenol, 2-[(1R)-1-[(1R)-1-phenylethyl]amino]ethyl)- (9CI)
 MF C16 H19 N O

Absolute stereochemistry. Rotation (+).



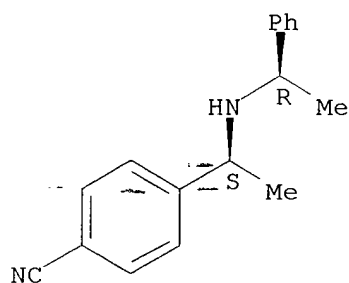
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Chromium, tricarbonyl[(1,2,3,4,5,6-eta.)-.alpha.-methyl-N-(1-phenylethyl)benzenemethanamine]-, [S-(R*,R*)]- (9CI)
 MF C19 H19 Cr N O3
 CI CCS



L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzonitrile, 4-[1-[(1-phenylethyl)amino]ethyl]-, (R*,S*)- (9CI)
 MF C17 H18 N2

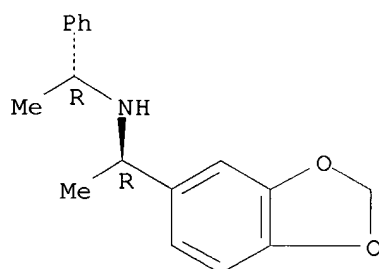
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 1,3-Benzodioxole-5-methanamine, .alpha.-methyl-N-[(1R)-1-phenylethyl]-,
 (.alpha.R)- (9CI)
 MF C17 H19 N O2

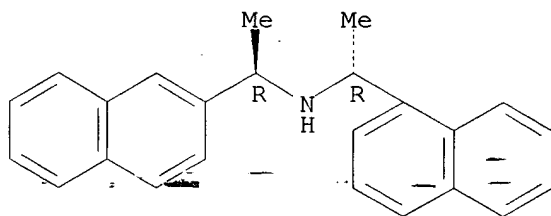
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 1-Naphthalenemethanamine,
 .alpha.-methyl-N-[(1R)-1-(2-naphthalenyl)ethyl]-,
 (.alpha.R)- (9CI)
 MF C24 H23 N
 CI COM

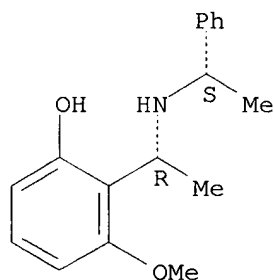
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Phenol, 3-methoxy-2-[1-[(1-phenylethyl)amino]ethyl]-, [S-(R*,S*)]- (9CI)
 MF C17 H21 N O2

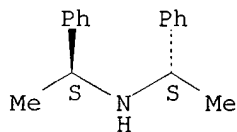
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzenemethanamine, .alpha.-methyl-N-[(1S)-1-phenylethyl]-, lithium salt, (.alpha.S)- (9CI)
 MF C16 H19 N . Li

Absolute stereochemistry. Rotation (-).

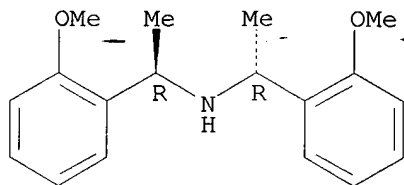


● Li

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzenemethanamine, 2-methoxy-N-[(1R)-1-(2-methoxyphenyl)ethyl]-.alpha.-methyl-, (.alpha.R)- (9CI)

MF C18 H23 N O2

Absolute stereochemistry.

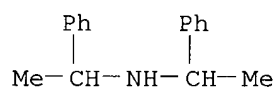


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN Benzenemethanamine, .alpha.-methyl-N-(1-phenylethyl)-, lithium salt (9CI)

MF C16 H19 N . Li



● Li

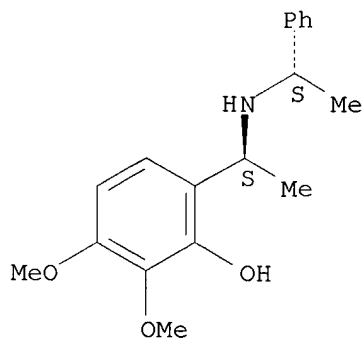
L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN Phenol, 2,3-dimethoxy-6-[1-[(1-phenylethyl)amino]ethyl]-, [S-(R*,R*)]-, perchlorate (salt) (9CI)

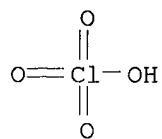
MF C18 H23 N O3 . Cl H O4

CM 1

Absolute stereochemistry.

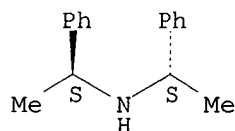


CM 2



L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzenemethanamine, .alpha.-methyl-N-[(1S)-1-phenylethyl]-,
 hydrochloride,
 (.alpha.S)- (9CI)
 MF C16 H19 N . Cl H

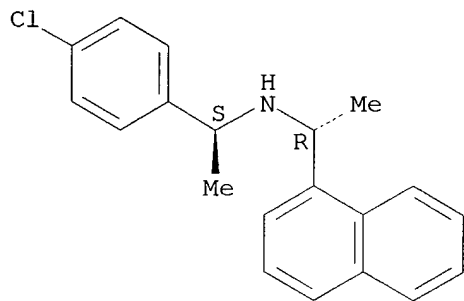
Absolute stereochemistry. Rotation (-).



● HCl

L5 252 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 1-Naphthalenemethanamine,
 N-[(1S)-1-(4-chlorophenyl)ethyl]-.alpha.-methyl-
 , (.alpha.R)- (9CI)
 MF C20 H20 Cl N

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

FULL ESTIMATED COST

ENTRY	SESSION
142.18	142.39

FILE 'CAPLUS' ENTERED AT 14:38:02 ON 24 APR 2002
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2002 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 24 Apr 2002 VOL 136 ISS 17
FILE LAST UPDATED: 23 Apr 2002 (20020423/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

CAS roles have been modified effective December 16, 2001. Please check your SDI profiles to see if they need to be revised. For information on CAS roles, enter HELP ROLES at an arrow prompt or use the CAS Roles thesaurus (/RL field) in this file.

=> l5

L6 233 L5

=> save temp l6 chiramines/a

ANSWER SET L6 HAS BEEN SAVED AS 'CHIRAMINES/A'

=> imine

 15447 IMINE
 10247 IMINES
L7 21166 IMINE
 (IMINE OR IMINES)

=> l6 and l7

L8 27 L6 AND L7

=> reduc?

 1633885 REDUC?
 723193 REDN
 37985 REDNS
 746911 REDN
 (REDN OR REDNS)
L9 2051132 REDUC?
 (REDUC? OR REDN)

=> l8 and l9

L10 11 L8 AND L9

=> d l10 1-11 ti

L10 ANSWER 1 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI Microwave-assisted catalytic intermolecular hydroamination of alkynes

L10 ANSWER 2 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI Asymmetric synthesis of planar chiral (arene)tricarbonylchromium complexes
 via enantioselective deprotonation by conformationally constrained chiral lithium-amide bases

L10 ANSWER 3 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI Asymmetric **reduction** of enantiopure **imines** with zinc borohydride: stereoselective synthesis of chiral amines

L10 ANSWER 4 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI Chemoselective **reductive** amination of aldehydes and ketones by dibutylchlorotin hydride-HMPA complex

L10 ANSWER 5 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI Aminoborohydrides. 6. Diastereoselective **reduction** of the carbon-nitrogen double bond in chiral **imines** using lithium diethylaminoborohydride and lithium diisopropylaminoborohydride

L10 ANSWER 6 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI Diastereoselective nucleophilic additions to **imines** attached to tricarbonyl(arene) chromium moieties

L10 ANSWER 7 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI NAD(P)+-NAD(P)H models. 70. **Reduction** of **imines** with Hantzsch ester in the presence of silica gel

L10 ANSWER 8 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI Synthesis of amines by **reduction** of **imines** with the MCl₂/NaBH₄ (M = cobalt, nickel) system

L10 ANSWER 9 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI Chiral building blocks for the synthesis of N-containing natural products.
 Part I. Enantiomerically pure oxygenated 1-phenylethylamines from substituted acetophenones: by **reductive** amination and regioselective benzylic cleavage

L10 ANSWER 10 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI NADH models. XXI. Stereoselective **reduction** of chiral **imines** with Hantzsch ester

L10 ANSWER 11 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI Electrosynthesis of N-phenylethyl-substituted DL-phenylglycine- and DL- α -phenylalanine esters via **reductive** carboxylation of N-benzylidene- and N-(α -methylbenzylidene)amines

=> save temp all reducamin/l

L# LIST L1-L10 HAS BEEN SAVED AS 'REDUCAMIN/L'

=> d l10 8 ti fbib abs

L10 ANSWER 8 OF 11 CAPLUS COPYRIGHT 2002 ACS
 TI Synthesis of amines by **reduction** of **imines** with the MCl₂/NaBH₄ (M = cobalt, nickel) system

AN 1989:533330 CAPLUS
 DN 111:133330
 TI Synthesis of amines by **reduction** of **imines** with the
 MCl₂/NaBH₄ (M = cobalt, nickel) system
 AU Periasamy, M.; Devasagayaraj, A.; Satyanarayana, N.; Narayana, C.
 CS Sch. Chem., Univ. Hyderabad, Hyderabad, 500 134, India
 SO Synth. Commun. (1989), 19(3-4), 565-73
 CODEN: SYNCAV; ISSN: 0039-7911
 DT Journal
 LA English
 OS CASREACT 111:133330
 AB **Imines**, anils, and enamines were **reduced** by CoCl₂ or
 NiCl₂/NaBH₄/CH₃OH in THF under mild conditions to afford the
 corresponding
 amines in 64-82% yields. Thus, PhCMe:NCHMePh was added to the
 CoCl₂-contg. reagent and the mixt. stirred 2 h at -10.degree. to give 64%
 (PhCHMe)₂NH.

=> logoff hold		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	12.23	154.62
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-0.62	-0.62

SESSION WILL BE HELD FOR 60 MINUTES
 STN INTERNATIONAL SESSION SUSPENDED AT 14:43:10 ON 24 APR 2002

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1623paz

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *
 SESSION RESUMED IN FILE 'CAPLUS' AT 14:44:49 ON 24 APR 2002
 FILE 'CAPLUS' ENTERED AT 14:44:49 ON 24 APR 2002
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COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	12.23	154.62
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-0.62	-0.62

=> d his

(FILE 'HOME' ENTERED AT 14:33:48 ON 24 APR 2002)

FILE 'REGISTRY' ENTERED AT 14:34:12 ON 24 APR 2002

L1 STRUCTURE UPLOADED

L2 6 SEARCH L1 SSS SAM
L3 STRUCTURE UPLOADED
L4 1 SEARCH L3 SSS SAM
L5 252 SEARCH L3 SSS FULL

FILE 'CAPLUS' ENTERED AT 14:38:02 ON 24 APR 2002

L6 233 L5
SAVE TEMP L6 CHIRAMINES/A
L7 21166 IMINE
L8 27 L6 AND L7
L9 2051132 REDUC?
L10 11 L8 AND L9
SAVE TEMP ALL REDUCAMIN/L

=> ?fluor?

L11 1097263 ?FLUOR?

=> l6 and l11

L12 22 L6 AND L11

=> l10 and l12

L13 1 L10 AND L12

=> d l13 ti fbib abs

L13 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2002 ACS

TI Asymmetric synthesis of planar chiral (arene)tricarbonylchromium
complexes

via enantioselective deprotonation by conformationally constrained chiral
lithium-amide bases

AN 2000:748505 CAPLUS

DN 134:42233

TI Asymmetric synthesis of planar chiral (arene)tricarbonylchromium
complexes

via enantioselective deprotonation by conformationally constrained chiral
lithium-amide bases

AU Pache, Sandrine; Botuha, Candice; Franz, Roberto; Kundig, E. Peter;
Einhorn, Jacques

CS Departement de Chimie Organique, Universite de Geneve, Geneva, CH-1211/4,
Switz.

SO Helvetica Chimica Acta (2000), 83(9), 2436-2451

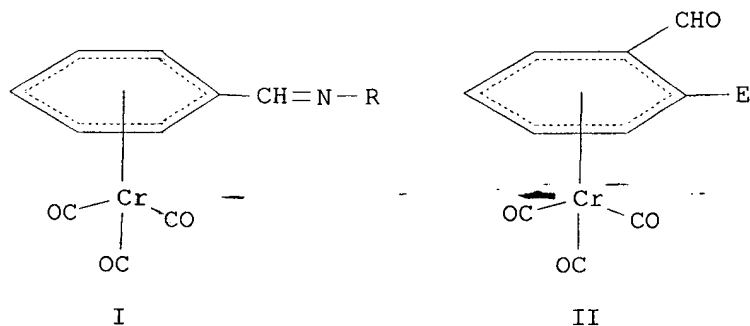
CODEN: HCACAV; ISSN: 0018-019X

PB Verlag Helvetica Chimica Acta

DT Journal

LA English

GI



AB Enantioselective lithiation/electrophilic addn. reactions with eight chiral Li-amide bases, 1-8, and five [Cr(arene)(CO)₃] complexes, 9-13, were studied. Restriction of conformational freedom in the chiral Li-amide base Li-1, in general, did not result in an increase in asym. induction. A new route to enantiomerically enriched (75-92%) planar chiral ortho-substituted benzaldehyde complexes via enantioselective lithiation of benzaldimine complexes 16 and 17 (I; R = cyclohexyl, Ph) is reported. Within the (1S)-enantiomer series of o-substituted benzaldehyde complexes 18a-d (II; E = Me₃Si, Me₃Sn, Me, COOMe), generated from 16 and 17, the sign of the sp. rotation, [α]_{D20}, is pos., except for the trimethylstannyl deriv. 18b. This is interpreted in terms of a reversed conformation of the aldehyde group.

RE.CNT 82 THERE ARE 82 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> l12 not l13

L14 21 L12 NOT L13

=> save temp l14 fluoroamine/a

ANSWER SET L14 HAS BEEN SAVED AS 'FLUOROAMINE/A'

=> d l14 1-21 ti

L14 ANSWER 1 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Preparation of optically active .alpha.-methylbis-3,5-(**trifluoromethyl**)benzylamine from optically active imines

L14 ANSWER 2 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Purification of optically active .alpha.-methyl-3,5-bis(**trifluoromethyl**)benzylamines

L14 ANSWER 3 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Enantioselective synthesis of cyclohexenylalkenes by asymmetric deprotonation of 4-tert-butylcyclohexanone followed by O-nonaflation and Heck couplings

L14 ANSWER 4 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Preparation of 1-arylethylamines as calcium receptor ligands

L14 ANSWER 5 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI 195Pt NMR determination of the enantiomeric purity and absolute configuration of trisubstituted allenes by using [PtCl₃(C₂H₄)]-[(S,S)-(1-

NpMeCH)2NH2]+ as CDA

L14 ANSWER 6 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Desymmetrization of 4,4-disubstituted cyclohexanones by enzyme-catalyzed resolution of their enol acetates

L14 ANSWER 7 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI A practical o-hydroxybenzylamines promoted enantioselective addition of dialkylzincs to aldehydes with asymmetric amplification

L14 ANSWER 8 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Aldol Addition of Lithium and Boron Enolates of 1,3-Dioxan-5-ones to Aldehydes. A New Entry into Monosaccharide Derivatives

L14 ANSWER 9 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI The asymmetric synthesis of phosphorus- and sulfur-containing tricarbonyl(.eta.6-arene)chromium complexes using the chiral base approach

L14 ANSWER 10 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Optical Resolution and Epimerization of **Fluorosilane** Having an Optically Active Amino Group: A New, Convenient Access to Optically Active Silicon Compounds

L14 ANSWER 11 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI New symmetry-breaking deprotonation reactions of cyclic imides using a chiral lithium amide base

L14 ANSWER 12 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Enantiomeric impurities in chiral catalysts, auxiliaries and synthons used in enantioselective synthesis

L14 ANSWER 13 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Chiral base mediated asymmetric synthesis of tricarbonyl(.eta.6-arene)chromium complexes

L14 ANSWER 14 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Diastereoselective addition of methyllithium and dimethylcuprate-boron **trifluoride** to imines derived from (S)-1-phenylethylamine

L14 ANSWER 15 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Cesium **fluoride**-mediated Horner-Wittig addition reactions of silyl phosphine oxides; synthesis of optically active silyl phosphine oxides using chiral bases and a chiral acid

L14 ANSWER 16 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Synthesis of .alpha.-**trifluoromethyl** substituted .alpha.-amino acid derivatives from methyl 3,3,3-**trifluoro**-2-diazopropionate

L14 ANSWER 17 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Preparation of substituted N-aryl-1,2-diaminocyclobutene-3,4-dione smooth muscle relaxants

L14 ANSWER 18 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Concerning the asymmetric metalation of ferrocenes by chiral lithium amide bases

L14 ANSWER 19 OF 21 CAPLUS COPYRIGHT 2002 ACS
TI Diastereoselective addition of methylcopper- and dimethylcuprate-boron
trifluoride reagents to (S)-N-alkylidene-1-phenylethylamines

L14 ANSWER 20 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Chiral building blocks for the synthesis of N-containing natural
products.

4. A facile method for the asymmetric synthesis of enantiomerically pure
1-(2-**fluorophenyl**)ethylamine

L14 ANSWER 21 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Enantiotopic interactions in the **fluorescence** quenching of
camphor by chiral amines

=> d l14 1-2 ti fbib abs

L14 ANSWER 1 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Preparation of optically active .alpha.-methylbis-3,5-(
trifluoromethyl)benzylamine from optically active imines

AN 2002:77456 CAPLUS

DN 136:134565

TI Preparation of optically active .alpha.-methylbis-3,5-(
trifluoromethyl)benzylamine from optically active imines

IN Ishii, Akio; Kuriyama, Suguru; Kanai, Masatomi; Hayami, Takashi

PA Central Glass Co., Ltd., Japan

SO Jpn. Kokai Tokkyo Koho, 9 pp.

CODEN: JKXXAF

DT Patent

LA Japanese

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 2002030048	A2	20020129	JP 2000-142460	20000515
				JP 2000-138349 A	20000511

OS CASREACT 136:134565; MARPAT 136:134565

AB (R)- or (S)-H₂NCHMeC₆H₃(CF₃)₂-3,5, useful as an intermediate for
pesticides and drugs, is prepd. by asym. hydrogenation of optically
active-ArCHMeN:CMcC₆H₃(CF₃)₂-3,5 (Ar = Ph, 1- or 2-naphthyl) and
hydrogenolysis of the resulting optically active ArCHMeNHCHMeC₆H₃(CF₃)₂-
3,5. Thus, MeCOC₆H₃(CF₃)₂-3,5 was refluxed with (S)-H₂NCHPhMe and
p-MeC₆H₄SO₃H to give quant. optically active imine, which was
hydrogenated

with NaBH₄ in EtOH to afford 7.8:1 (SS)- and (SR)-PhCHMeNHCHMeC₆H₃(CF₃)₂-
3,5 with 100% conversion. The diastereomeric mixt. was hydrogenated over
Pd/C to give 75% (S)-H₂NCHMeC₆H₃(CF₃)₂-3,5 with 76% ee.

L14 ANSWER 2 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Purification of optically active .alpha.-methyl-3,5-bis(
trifluoromethyl)benzylamines

AN 2002:23507 CAPLUS

DN 136:69641

TI Purification of optically active .alpha.-methyl-3,5-bis(
trifluoromethyl)benzylamines

IN Ishii, Akio; Kuriyama, Masaru; Yasumoto, Manabu; Kanai, Masatomi; Hayami,
Takashi

PA Central Glass Co., Ltd., Japan

SO Jpn. Kokai Tokkyo Koho, 8 pp.

CODEN: JKXXAF

~~LE~~ Patent
~~LA~~ Japanese
FIN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 2002003453	A2	20020109	JP 2000-185405	20000620
CS	MARPAT 136:69641				
AB	The compds. 3,5-(F3C)2C6H3CHMeNHR (R = H, benzyl, aryl, CHMeAr; Ar = Ph, 1- or 2-naphthyl) are purified by converting into inorg. acid or org. acid salts and recrystn. .alpha.-Methyl-bis-3,5-(trifluoromethyl)benzylamine (S-isomer:R-isomer = 7.4:1) was reacted with p-MeC6H4SO3H in PhMe at 60-70.degree. for 30 min to give (S)-.alpha.-methyl-3,5-bis(trifluoromethyl)benzylamine p-toluenesulfonate with 82.7% e.e.				

=> logoff hold

CST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
29.45	171.84

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
-2.48	-2.48

CA SUBSCRIBER PRICE

SESSION WILL BE HELD FOR 60 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 14:49:03 ON 24 APR 2002

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NEWS 1		Web Page URLs for STN Seminar Schedule - N. America
NEWS 2	Jan 25	BLAST(R) searching in REGISTRY available in STN on the Web
NEWS 3	Jan 29	FSTA has been reloaded and moves to weekly updates
NEWS 4	Feb 01	DKILIT now produced by FIZ Karlsruhe and has a new update frequency
NEWS 5	Feb 19	Access via Tymnet and SprintNet Eliminated Effective 3/31/02
NEWS 6	Mar 08	Gene Names now available in BIOSIS
NEWS 7	Mar 22	TOXLIT no longer available
NEWS 8	Mar 22	TRCTHERMO no longer available
NEWS 9	Mar 28	US Provisional Priorities searched with P in CA/CAplus and USPATFULL
NEWS 10	Mar 28	LIPINSKI/CALC added for property searching in REGISTRY
NEWS 11	Apr 02	PAPERCHEM no longer available on STN. Use PAPERCHEM2 instead.
NEWS 12	Apr 08	"Ask CAS" for self-help around the clock
NEWS 13	Apr 09	BEILSTEIN: Reload and Implementation of a New Subject Area

NEWS 14 Apr 09 ZDB will be removed from STN
 NEWS 15 Apr 19 US Patent Applications available in IFICDB, IFIPAT, and
 IFIUDB
 NEWS 16 Apr 22 Records from IP.com available in CAPLUS, HCAPLUS, and
 ZCAPLUS
 NEWS 17 Apr 22 BIOSIS Gene Names now available in TOXCENTER
 NEWS 18 Apr 22 Federal Research in Progress (FEDRIP) now available

NEWS EXPRESS February 1 CURRENT WINDOWS VERSION IS V6.0d,
 CURRENT MACINTOSH VERSION IS V6.0a(ENG) AND V6.0Ja(JP),
 AND CURRENT DISCOVER FILE IS DATED 05 FEBRUARY 2002
 NEWS HOURS STN Operating Hours Plus Help Desk Availability
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 NEWS WWW CAS World Wide Web Site (general information)

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FILE CONTAINS CURRENT INFORMATION.

LAST RELOADED: Apr 19, 2002 (20020419/UP).

=>

NAME	CREATED	NOTES/TITLE
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AIDSMOTIV/A	TEMP	117 ANSWERS IN FILE CAPLUS
AIDSSRCH/L	TEMP	5 L-NUMBERS
ALKYLATIN/L	13 DEC 2001	9 L-NUMBERS
ANDROELECTRO/A	TEMP	10 ANSWERS IN FILE CAPLUS
ANDRONUCLEO/A	TEMP	4173 ANSWERS IN FILE CAPLUS
ANDROSRCH/L	TEMP	23 L-NUMBERS
ANTI HIV/A	TEMP	6 ANSWERS IN FILE CAPLUS
ANTI HIVFREE/A	TEMP	41 ANSWERS IN FILE CAPLUS
CHIRAMINES/A	TEMP	233 ANSWERS IN FILE CAPLUS
FLUOROAMINE/A	TEMP	21 ANSWERS IN FILE CAPLUS

HIVCMPDS/A	TEMP	10 ANSWERS IN FILE CAPLUS
INDIUMCL3/A	30 MAY 2001	1 ANSWER IN FILE REGISTRY
LTWENTAUGFOR/A	04 AUG 2001	72 ANSWERS IN FILE CAPLUS
NEOTAMECRYST/A	24 APR 2001	59 ANSWERS IN FILE CAPLUS
NSAIDTRGT/A	TEMP	10 ANSWERS IN FILE CAPLUS
NVLARMFULGEN/A	19 APR 2001	196 ANSWERS IN FILE REGISTRY
POHBENZALDEH/A	10 JUL 2001	5519 ANSWERS IN FILE CAPLUS
PROSTACMPD15/A	01 AUG 2001	34 ANSWERS IN FILE CAPLUS
RANEYSRCH/L	TEMP	16 L-NUMBERS
REDUCAMIN/L	TEMP	10 L-NUMBERS
STILLEAPP/L	07 JAN 2002	17 L-NUMBERS
TWOAMINOPOLY/Q	16 APR 2001	UPLOADED STRUCTURE

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COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.12	0.33

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FILE COVERS 1907 - 25 Apr 2002 VOL 136 ISS 17
 FILE LAST UPDATED: 23 Apr 2002 (20020423/ED)

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CAS roles have been modified effective December 16, 2001. Please check your SDI profiles to see if they need to be revised. For information on CAS roles, enter HELP ROLES at an arrow prompt or use the CAS Roles thesaurus (/RL field) in this file.

=>

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L1          STR
L2  (       252)SEA FILE=REGISTRY SSS FUL L1
L3          233 SEA FILE=CAPLUS ABB=ON  PLU=ON  L2

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=>

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L4          STR
L5  (       252)SEA FILE=REGISTRY SSS FUL L4
L6  (       233)SEA FILE=CAPLUS ABB=ON  PLU=ON  L5

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L7 (21166)SEA FILE=CAPLUS ABB=ON PLU=ON IMINE
 L8 (27)SEA FILE=CAPLUS ABB=ON PLU=ON L6 AND L7
 L9 (2051132)SEA FILE=CAPLUS ABB=ON PLU=ON REDUC?
 L10 (11)SEA FILE=CAPLUS ABB=ON PLU=ON L8 AND L9
 L11 (1097263)SEA FILE=CAPLUS ABB=ON PLU=ON ?FLUOR?
 L12 (22)SEA FILE=CAPLUS ABB=ON PLU=ON L6 AND L11
 L13 (1)SEA FILE=CAPLUS ABB=ON PLU=ON L10 AND L12
 L14 (21)SEA FILE=CAPLUS ABB=ON PLU=ON L12 NOT L13

=>

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.40	0.73

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FILE CONTAINS CURRENT INFORMATION.
 LAST RELOADED: Apr 19, 2002 (20020419/UP).

=>

L15 STR
 L16 (6)SEA FILE=REGISTRY SSS SAM L15
 L17 STR
 L18 (1)SEA FILE=REGISTRY SSS SAM L17
 L19 (252)SEA FILE=REGISTRY SSS FUL L17
 L20 (233)SEA FILE=CAPLUS ABB=ON PLU=ON L19
 L21 (21166)SEA FILE=CAPLUS ABB=ON PLU=ON IMINE
 L22 (27)SEA FILE=CAPLUS ABB=ON PLU=ON L20 AND L21
 L23 (2051132)SEA FILE=CAPLUS ABB=ON PLU=ON REDUC?
 L24 (11)SEA FILE=CAPLUS ABB=ON PLU=ON L22 AND L23

=> logoff hold

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.06	0.79

SESSION WILL BE HELD FOR 60 MINUTES
 STN INTERNATIONAL SESSION SUSPENDED AT 07:06:39 ON 25 APR 2002

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LOGINID:sssptal623paz

PASSWORD:

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COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
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FULL ESTIMATED COST	0.06	0.79
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=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.06	0.79

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 DICTIONARY FILE UPDATES: 23 APR 2002 HIGHEST RN 406909-40-8

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Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES
 for more information. See STN Note 27, Searching Properties in the CAS
 Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=> file caplus		
COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.38	1.17

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FILE COVERS 1907 - 25 Apr 2002 VOL 136 ISS 17
 FILE LAST UPDATED: 23 Apr 2002 (20020423/ED)

This file contains CAS Registry Numbers for easy and accurate
 substance identification.

CAS roles have been modified effective December 16, 2001. Please
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 information on CAS roles, enter HELP ROLES at an arrow prompt or use
 the CAS Roles thesaurus (/RL field) in this file.

=> d his

(FILE 'HOME' ENTERED AT 07:04:49 ON 25 APR 2002)

FILE 'STNGUIDE' ENTERED AT 07:05:00 ON 25 APR 2002

FILE 'CAPLUS' ENTERED AT 07:05:57 ON 25 APR 2002

ACT CHIRAMINES/A

L1 STR
L2 (252)SEA FILE=REGISTRY SSS FUL L1
L3 233 SEA FILE=CAPLUS ABB=ON PLU=ON L2

ACT FLUOROAMINE/A

L4 STR
L5 (252)SEA FILE=REGISTRY SSS FUL L4
L6 (233)SEA FILE=CAPLUS ABB=ON PLU=ON L5
L7 (21166)SEA FILE=CAPLUS ABB=ON PLU=ON IMINE
L8 (27)SEA FILE=CAPLUS ABB=ON PLU=ON L6 AND L7
L9 (2051132)SEA FILE=CAPLUS ABB=ON PLU=ON REDUC?
L10 (11)SEA FILE=CAPLUS ABB=ON PLU=ON L8 AND L9
L11 (1097263)SEA FILE=CAPLUS ABB=ON PLU=ON ?FLUOR?
L12 (22)SEA FILE=CAPLUS ABB=ON PLU=ON L6 AND L11
L13 (1)SEA FILE=CAPLUS ABB=ON PLU=ON L10 AND L12
L14 21 SEA FILE=CAPLUS ABB=ON PLU=ON L12 NOT L13

FILE 'STNGUIDE' ENTERED AT 07:06:00 ON 25 APR 2002

ACT REDUCAMIN/L

L15 STR
L16 (6)SEA FILE=REGISTRY SSS SAM L15
L17 STR
L18 (1)SEA FILE=REGISTRY SSS SAM L17
L19 (252)SEA FILE=REGISTRY SSS FUL L17
L20 (233)SEA FILE=CAPLUS ABB=ON PLU=ON L19
L21 (21166)SEA FILE=CAPLUS ABB=ON PLU=ON IMINE
L22 (27)SEA FILE=CAPLUS ABB=ON PLU=ON L20 AND L21
L23 (2051132)SEA FILE=CAPLUS ABB=ON PLU=ON REDUC?
L24 (11)SEA FILE=CAPLUS ABB=ON PLU=ON L22 AND L23

FILE 'REGISTRY' ENTERED AT 07:15:34 ON 25 APR 2002

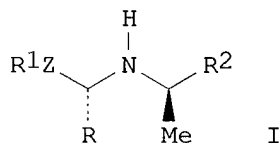
FILE 'CAPLUS' ENTERED AT 07:15:40 ON 25 APR 2002

=> d l14 4,20 ti fbib abs

L14 ANSWER 4 OF 21 CAPLUS COPYRIGHT 2002 ACS
TI Preparation of 1-arylethylamines as calcium receptor ligands
AN 2001:241760 CAPLUS
DN 134:280612
TI Preparation of 1-arylethylamines as calcium receptor ligands
IN Van Wagenen, Bradford C.; Moe, Scott T.; Balandrin, Manuel F.; Delmar,
Eric G.; Nemeth, Edward F.
PA NPS Pharmaceuticals, Inc., USA
SO U.S., 142 pp.
CODEN: USXXAM

DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 6211244	B1	20010403	US 1995-546998	19951023
OS	MARPAT 134:280612				
GI					



AB Title compds., e.g., I [R = H or alkyl; R1,R2 = (un)substituted Ph or naphthyl; Z = (CH2)0-3] were prepd. Thus, (R)-1-(1-naphthyl)ethylamine was condensed with 2-acetonaphthone to give I (R = Me, R1 = 2-naphthyl,

R2 = 1-naphthyl, Z = bond). Data for biol. activity of title compds. were given.

RE.CNT 229 THERE ARE 229 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L14 ANSWER 20 OF 21 CAPLUS COPYRIGHT 2002 ACS

TI Chiral building blocks for the synthesis of N-containing natural products.

4. A facile method for the asymmetric synthesis of enantiomerically pure 1-(2-**fluorophenyl**)ethylamine

AN 1990:630871 CAPLUS

DN 113:230871

TI Chiral building blocks for the synthesis of N-containing natural products.

4. A facile method for the asymmetric synthesis of enantiomerically pure 1-(2-**fluorophenyl**)ethylamine

AU Bringmann, G.; Geisler, J. P.

CS Inst. Org. Chem., Univ. Wuerzburg, Wuerzburg, D-8700, Fed. Rep. Ger.

SO J. Fluorine Chem. (1990), 49(1), 67-73

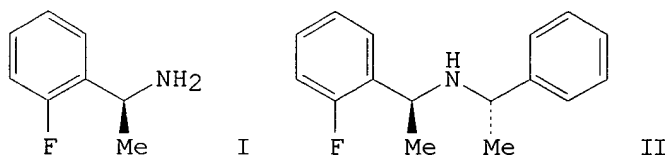
CODEN: JFLCAR; ISSN: 0022-1139

DT Journal

LA English

OS CASREACT 113:230871

GI



AB A simple, 2-step-procedure for the synthesis of optically active (S)-1-(2-

fluorophenyl)ethylamine (I) is described. Starting from com.

available 2-**fluoroacetophenone**, imination with
 (S)-1-phenylethylamine, followed by stereoselective hydrogenation over
 Raney Ni gives the secondary amine II. Subsequent regioselective
 hydrogenolytic cleavage of homogeneous II yields enantiomerically pure
 title compd.

```
=> logoff hold
COST IN U.S. DOLLARS          SINCE FILE      TOTAL
                                ENTRY      SESSION
FULL ESTIMATED COST          6.16          7.33

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)  SINCE FILE      TOTAL
                                                ENTRY      SESSION
CA SUBSCRIBER PRICE          -1.24         -1.24
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SESSION WILL BE HELD FOR 60 MINUTES
 STN INTERNATIONAL SESSION SUSPENDED AT 07:18:14 ON 25 APR 2002

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal623paz

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *
 SESSION RESUMED IN FILE 'CAPLUS' AT 08:08:32 ON 25 APR 2002
 FILE 'CAPLUS' ENTERED AT 08:08:32 ON 25 APR 2002
 COPYRIGHT (C) 2002 AMERICAN CHEMICAL SOCIETY (ACS)

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COST IN U.S. DOLLARS          SINCE FILE      TOTAL
                                ENTRY      SESSION
FULL ESTIMATED COST          6.16          7.33

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)  SINCE FILE      TOTAL
                                                ENTRY      SESSION
CA SUBSCRIBER PRICE          -1.24         -1.24
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=> d his

(FILE 'HOME' ENTERED AT 07:04:49 ON 25 APR 2002)

FILE 'STNGUIDE' ENTERED AT 07:05:00 ON 25 APR 2002

FILE 'CAPLUS' ENTERED AT 07:05:57 ON 25 APR 2002

ACT CHIRAMINES/A

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L1          STR
L2 (        252)SEA FILE=REGISTRY SSS FUL L1
L3          233 SEA FILE=CAPLUS ABB=ON  PLU=ON  L2
```

 ACT FLUOROAMINE/A

```
-----
L4          STR
L5 (        252)SEA FILE=REGISTRY SSS FUL L4
L6 (        233)SEA FILE=CAPLUS ABB=ON  PLU=ON  L5
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```

L7 (      21166)SEA FILE=CAPLUS ABB=ON  PLU=ON  IMINE
L8 (      27)SEA FILE=CAPLUS ABB=ON  PLU=ON  L6 AND L7
L9 (    2051132)SEA FILE=CAPLUS ABB=ON  PLU=ON  REDUC?
L10 (      11)SEA FILE=CAPLUS ABB=ON  PLU=ON  L8 AND L9
L11 (   1097263)SEA FILE=CAPLUS ABB=ON  PLU=ON  ?FLUOR?
L12 (      22)SEA FILE=CAPLUS ABB=ON  PLU=ON  L6 AND L11
L13 (      1)SEA FILE=CAPLUS ABB=ON  PLU=ON  L10 AND L12
L14 (     21)SEA FILE=CAPLUS ABB=ON  PLU=ON  L12 NOT L13
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FILE 'STNGUIDE' ENTERED AT 07:06:00 ON 25 APR 2002
 ACT REDUCAMIN/L

```

L15          STR
L16 (      6)SEA FILE=REGISTRY SSS SAM L15
L17          STR
L18 (      1)SEA FILE=REGISTRY SSS SAM L17
L19 (     252)SEA FILE=REGISTRY SSS FUL L17
L20 (     233)SEA FILE=CAPLUS ABB=ON  PLU=ON  L19
L21 (   21166)SEA FILE=CAPLUS ABB=ON  PLU=ON  IMINE
L22 (      27)SEA FILE=CAPLUS ABB=ON  PLU=ON  L20 AND L21
L23 (   2051132)SEA FILE=CAPLUS ABB=ON  PLU=ON  REDUC?
L24 (      11)SEA FILE=CAPLUS ABB=ON  PLU=ON  L22 AND L23
-----

```

FILE 'REGISTRY' ENTERED AT 07:15:34 ON 25 APR 2002

FILE 'CAPLUS' ENTERED AT 07:15:40 ON 25 APR 2002

=> mandelic

L25 4560 MANDELIC

=> tartaric

28287 TARTARIC

1 TARTARICS

L26 28288 TARTARIC

(TARTARIC OR TARTARICS)

=> 125 or 126

L27 32468 L25 OR L26

=> 13 and 127

L28 8 L3 AND L27

=> d 128 1-8 ti

L28 ANSWER 1 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Purification of optically active .alpha.-methyl-3,5-bis(trifluoromethyl)benzylamines

L28 ANSWER 2 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Preparation of optically active .alpha.-methylbenzylamine

L28 ANSWER 3 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Design, synthesis, and optical resolution of a novel non-natural chiral auxiliary, 1-(2,5-dimethoxyphenyl)ethylamine. Application to diastereoselective alkylation of aldimines

L28 ANSWER 4 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Preparation of 6-aryl-(methyl- or methylenedene)-quinoline derivatives as voltage-gated potassium channel blockers

L28 ANSWER 5 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Preparation of optically active .alpha.-methylbenzylamine

L28 ANSWER 6 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Habit modification of a diastereomeric salt with an additive in optical resolution

L28 ANSWER 7 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Chiral base-induced [2,3] Wittig rearrangement of acyclic .alpha.-(propargyloxy)acetic acids and amides

L28 ANSWER 8 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Synthesis of stereomeric .alpha.,.alpha.'-dimethyldibenzylamines and their benzoates

=> d 128 1,2,5,6,8 ti fbib abs

L28 ANSWER 1 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Purification of optically active .alpha.-methyl-3,5-bis(trifluoromethyl)benzylamines

AN 2002:23507 CAPLUS

DN 136:69641

TI Purification of optically active .alpha.-methyl-3,5-bis(trifluoromethyl)benzylamines

IN Ishii, Akio; Kuriyama, Masaru; Yasumoto, Manabu; Kanai, Masatomi; Hayami, Takashi

PA Central Glass Co., Ltd., Japan

SO Jpn. Kokai Tokkyo Koho, 8 pp.

CODEN: JKXXAF

DT Patent

LA Japanese

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 2002003453	A2	20020109	JP 2000-185405	20000620
OS	MARPAT 136:69641				
AB	The compds. 3,5-(F3C)2C6H3CHMeNHR (R = H, benzyl, aryl, CHMeAr; Ar = Ph, 1- or 2-naphthyl) are purified by converting into inorg. acid or org. acid				

salts and recrystn. .alpha.-Methyl-bis-3,5-(trifluoromethyl)benzylamine (S-isomer:R-isomer = 7.4:1) was reacted with p-MeC6H4SO3H in PhMe at 60-70.degree. for 30 min to give (S)-.alpha.-methyl-3,5-bis(trifluoromethyl)benzylamine p-toluenesulfonate with 82.7% e.e.

L28 ANSWER 2 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Preparation of optically active .alpha.-methylbenzylamine

AN 2000:750318 CAPLUS

DN 133:296269

TI Preparation of optically active .alpha.-methylbenzylamine

IN Murakami, Naomichi; Sakai, Kenichi; Tobiyama, Tadashi

PA Yamakawa Chemical Industry Co., Ltd., Japan

SO Jpn. Kokai Tokkyo Koho, 9 pp.

CODEN: JKXXAF

DT Patent

LA Japanese

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 2000297066	A2	20001024	JP 1999-104499	19990412
AB	Title compd. (I) is prepd. by treatment of (RS)-I with optically active mandelic acid in aq. media and optical resoln. of diastereomer salt, wherein at least a part of the (RS)-I is prepd. by racemization of optically active I in the presence of catalytic amt. of strongly basic racemization agents and contains bis(.alpha.-methylbenzyl)amine (II). A mixt. of 163.5 kg (RS)-I and 214 kg crude (RS)-I (contg. 1.5 kg II) was treated with Na (R)-mandelate in H2O in the presence of HCl at 5.degree. for .apprx.1 h to give diastereomer salt, which was decompd. with NaOH to give 64.6% (R)-I with 98.6% ee. (S)-I recovered from the process was racemized with NaH and used as a starting material.				

L28 ANSWER 5 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Preparation of optically active .alpha.-methylbenzylamine

AN 1994:298229 CAPLUS

DN 120:298229

TI Preparation of optically active .alpha.-methylbenzylamine

IN Sakai, Kenichi; Murakami, Naomichi; Saigo, Kazuhiko; Nohira, Hiroyuki

PA Yamakawa Chemical Ind, Japan

SO Jpn. Kokai Tokkyo Koho, 6 pp.

CODEN: JKXXAF

DT Patent

LA Japanese

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 06001757	A2	19940111	JP 1992-158356	19920617
	JP 3178086	B2	20010618		
AB	In the prepn. of (R)- or (S)-.alpha.-methylbenzylamine (I) from (RS)-I by diastereomeric method, (RS)-I is treated with optically active mandelic acid (II) in water solvents in the presence of bis(.alpha.-methylbenzyl)amine (III), and (in)org. acid salts of NH3 or primary amines and/or water-sol. inorg. salts, then optically active I.II obtained is crystd. as filterable crystals. Aq. soln. of 10 g (RS)-I was treated with 6.91 g (R)-II and aq. HCl, mixed with (R,R)-III.HCl under heating, then cooled to give 7.65 g (R)-I.II contg. (R)-I of 98.6% e.e.				

L28 ANSWER 6 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Habit modification of a diastereomeric salt with an additive in optical resolution

AN 1992:623398 CAPLUS

DN 117:223398

TI Habit modification of a diastereomeric salt with an additive in optical resolution

AU Sakai, Kenichi; Maekawa, Yasunari; Saigo, Kazuhiko; Sukegawa, Makoto; Murakami, Hisamichi; Nohira, Hiroyuki

CS Fac. Eng., Saitama Univ., Urawa, 338, Japan

SO Bull. Chem. Soc. (1992), 65(7), 1747-50

CODEN: BCSJA8; ISSN: 0009-2673

DT Journal

LA English

AB In the optical resoln. of .alpha.-methylbenzylamine (I) with **mandelic** acid (II), the dimeric deriv. of I, bis(.alpha.-methylbenzyl)amine (IV), caused a habit modification of the diastereomeric

salt, (R)-I.cntdot.(R)-II (III). The habit modification was strongly influenced by the stereochem. of IV. Amine (R,R)-IV changed the morphol. of the crystal of the diastereomeric salt, even at a concn. of 0.007 mol% of III; the shape of the crystal became a hexagonal plate from a long hexagonal plate, whereas its stereoisomer ((R,S)-IV) changed the morphol. less than (R,R)-IV and (S,S)-IV did not change at all. The habit modification of III by IV is discussed on the basis of the crystal growth mechanism while considering the stereochem. of IV.

L28 ANSWER 8 OF 8 CAPLUS COPYRIGHT 2002 ACS

TI Synthesis of stereomeric .alpha.,.alpha.'-dimethyldibenzylamines and their

benzoates

AN 1987:575567 CAPLUS

DN 107:175567

TI Synthesis of stereomeric .alpha.,.alpha.'-dimethyldibenzylamines and their

benzoates

AU Jablonska-Pikus, Teresa; Kurys, Krystyna; Janczewski, Marian

CS Inst. Chem., Uniw. M. Curie-Sklodowskiej, Lublin, Pol.

SO Ann. Univ. Mariae Curie-Sklodowska, Sect. AA: Phys. Chem. (1983), 38 119-29

CODEN: ACFCAD; ISSN: 0365-1193

DT Journal

LA Polish

AB The meso form and racemate were isolated from a mixt. of the stereoisomeric title amines. The latter was resolved into enantiomers by crystn. with (+)-tartaric acid. The benzoates of the enantiomers also were prepd.

=> logoff hold

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

26.69

27.86

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

CA SUBSCRIBER PRICE

-4.34

-4.34

SESSION WILL BE HELD FOR 60 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 08:13:41 ON 25 APR 2002

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal623paz

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *

SESSION RESUMED IN FILE 'CAPLUS' AT 08:26:43 ON 25 APR 2002

FILE 'CAPLUS' ENTERED AT 08:26:43 ON 25 APR 2002

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COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST	26.69	27.86
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-4.34	-4.34

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COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	26.69	27.86

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-4.34	-4.34

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STRUCTURE FILE UPDATES: 23 APR 2002 HIGHEST RN 406909-40-8
 DICTIONARY FILE UPDATES: 23 APR 2002 HIGHEST RN 406909-40-8

TSCA INFORMATION NOW CURRENT THROUGH July 7, 2001

Please note that search-term pricing does apply when
 conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES
 for more information. See STNote 27, Searching Properties in the CAS
 Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

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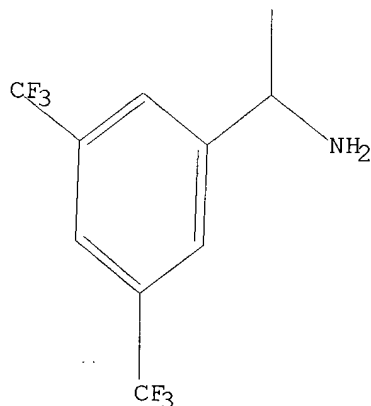
Uploading 09853085 bistrifluoromethylamine.str

L29 STRUCTURE UPLOADED

=> d 129

L29 HAS NO ANSWERS

L29 STR



Structure attributes must be viewed using STN Express query preparation.

=> search l29 exact full

FULL SEARCH INITIATED 08:28:31 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 14 TO ITERATE

100.0% PROCESSED 14 ITERATIONS

3 ANSWERS

SEARCH TIME: 00.00.01

L30 3 SEA EXA FUL L29

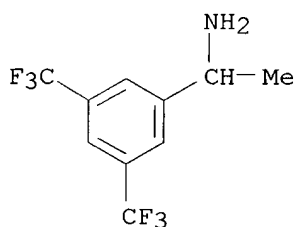
=> d scan

L30 3 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN Benzenemethanamine, .alpha.-methyl-3,5-bis(trifluoromethyl)- (9CI)

MF C10 H9 F6 N

CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):3

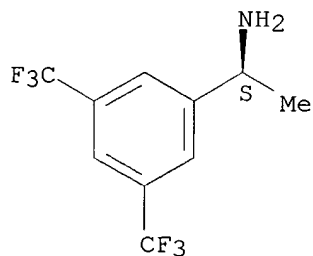
L30 3 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN Benzenemethanamine, .alpha.-methyl-3,5-bis(trifluoromethyl)-, (.alpha.S)- (9CI)

MF C10 H9 F6 N

CI COM

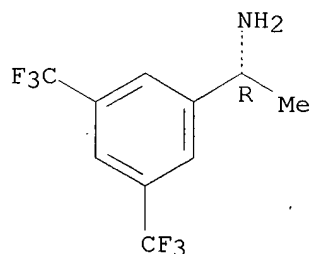
Absolute stereochemistry. Rotation (-).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L30 3 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenemethanamine, .alpha.-methyl-3,5-bis(trifluoromethyl)-, (.alpha.R)-
(9CI)
MF C10 H9 F6 N
CI COM

Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
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FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
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CA SUBSCRIBER PRICE

FILE 'CAPLUS' ENTERED AT 08:28:56 ON 25 APR 2002

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FILE COVERS 1907 - 25 Apr 2002 VOL 136 ISS 17

FILE LAST UPDATED: 23 Apr 2002 (20020423/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

CAS roles have been modified effective December 16, 2001. Please check your SDI profiles to see if they need to be revised. For information on CAS roles, enter HELP ROLES at an arrow prompt or use the CAS Roles thesaurus (/RL field) in this file.

=> 130

L31 12 L30

=> d his

(FILE 'HOME' ENTERED AT 07:04:49 ON 25 APR 2002)

FILE 'STNGUIDE' ENTERED AT 07:05:00 ON 25 APR 2002

FILE 'CAPLUS' ENTERED AT 07:05:57 ON 25 APR 2002

ACT CHIRAMINES/A

L1 STR
L2 (252)SEA FILE=REGISTRY SSS FUL L1
L3 233 SEA FILE=CAPLUS ABB=ON PLU=ON L2

ACT FLUOROAMINE/A

L4 STR
L5 (252)SEA FILE=REGISTRY SSS FUL L4
L6 (233)SEA FILE=CAPLUS ABB=ON PLU=ON L5
L7 (21166)SEA FILE=CAPLUS ABB=ON PLU=ON IMINE
L8 (27)SEA FILE=CAPLUS ABB=ON PLU=ON L6 AND L7
L9 (2051132)SEA FILE=CAPLUS ABB=ON PLU=ON REDUC?
L10 (11)SEA FILE=CAPLUS ABB=ON PLU=ON L8 AND L9
L11 (1097263)SEA FILE=CAPLUS ABB=ON PLU=ON ?FLUOR?
L12 (22)SEA FILE=CAPLUS ABB=ON PLU=ON L6 AND L11
L13 (1)SEA FILE=CAPLUS ABB=ON PLU=ON L10 AND L12
L14 21 SEA FILE=CAPLUS ABB=ON PLU=ON L12 NOT L13

FILE 'STNGUIDE' ENTERED AT 07:06:00 ON 25 APR 2002

ACT REDUCAMIN/L

L15 STR
L16 (6)SEA FILE=REGISTRY SSS SAM L15
L17 STR
L18 (1)SEA FILE=REGISTRY SSS SAM L17
L19 (252)SEA FILE=REGISTRY SSS FUL L17
L20 (233)SEA FILE=CAPLUS ABB=ON PLU=ON L19
L21 (21166)SEA FILE=CAPLUS ABB=ON PLU=ON IMINE
L22 (27)SEA FILE=CAPLUS ABB=ON PLU=ON L20 AND L21
L23 (2051132)SEA FILE=CAPLUS ABB=ON PLU=ON REDUC?
L24 (11)SEA FILE=CAPLUS ABB=ON PLU=ON L22 AND L23

FILE 'REGISTRY' ENTERED AT 07:15:34 ON 25 APR 2002

FILE 'CAPLUS' ENTERED AT 07:15:40 ON 25 APR 2002

L25 4560 MANDELIC
L26 28288 TARTARIC
L27 32468 L25 OR L26
L28 8 L3 AND L27

FILE 'REGISTRY' ENTERED AT 08:26:58 ON 25 APR 2002

L29 STRUCTURE UPLOADED
L30 3 SEARCH L29 EXACT FULL

FILE 'CAPLUS' ENTERED AT 08:28:56 ON 25 APR 2002

L31 12 L30

=> l31 and l27

L32 1 L31 AND L27

=> d l32 ti fbib abs

L32 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2002 ACS
TI Purification of optically active .alpha.-methyl-3,5-
bis(trifluoromethyl)benzylamines
AN 2002:23507 CAPLUS
DN 136:69641
TI Purification of optically active .alpha.-methyl-3,5-
bis(trifluoromethyl)benzylamines
IN Ishii, Akio; Kuriyama, Masaru; Yasumoto, Manabu; Kanai, Masatomi; Hayami,
Takashi
PA Central Glass Co., Ltd., Japan
SO Jpn. Kokai Tokkyo Koho, 8 pp.
CODEN: JKXXAF
DT Patent
LA Japanese
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 2002003453	A2	20020109	JP 2000-185405	20000620

OS MARPAT 136:69641
AB The compds. 3,5-(F3C)2C6H3CHMeNHR (R = H, benzyl, aryl, CHMeAr; Ar = Ph,
1- or 2-naphthyl) are purified by converting into inorg. acid or org.
acid
salts and recrystn. .alpha.-Methyl-bis-3,5-(trifluoromethyl)benzylamine
(S-isomer:R-isomer = 7.4:1) was reacted with p-MeC6H4SO3H in PhMe at
60-70.degree. for 30 min to give (S)-.alpha.-methyl-3,5-
bis(trifluoromethyl)benzylamine p-toluenesulfonate with 82.7% e.e.

=> d l31 1-12 ti

L31 ANSWER 1 OF 12 CAPLUS COPYRIGHT 2002 ACS
TI Preparation of optically active .alpha.-methylbis-3,5-
(trifluoromethyl)benzylamine from optically active imines

L31 ANSWER 2 OF 12 CAPLUS COPYRIGHT 2002 ACS
TI Purification of optically active .alpha.-methyl-3,5-
bis(trifluoromethyl)benzylamines

L31 ANSWER 3 OF 12 CAPLUS COPYRIGHT 2002 ACS
TI Preparation of cyclohexane derivatives for therapeutic use in the
treatment of disorders, such as depression, anxiety, pain, inflammation,
migraine, and vomiting

L31 ANSWER 4 OF 12 CAPLUS COPYRIGHT 2002 ACS
 TI Preparation of optically active .alpha.-methyl-bis-3,5-(trifluoromethyl)benzylamine

L31 ANSWER 5 OF 12 CAPLUS COPYRIGHT 2002 ACS
 TI Heteroaryl-containing thiourea derivatives useful as inhibitors of herpes viruses

L31 ANSWER 6 OF 12 CAPLUS COPYRIGHT 2002 ACS
 TI Aryl- and heteroaryl-substituted thiourea derivatives useful as inhibitors of herpes viruses

L31 ANSWER 7 OF 12 CAPLUS COPYRIGHT 2002 ACS
 TI Alpha-methylbenzyl-containing thiourea derivatives containing a phenylenediamine group, useful as inhibitors of herpes viruses

L31 ANSWER 8 OF 12 CAPLUS COPYRIGHT 2002 ACS
 TI Heterocyclic carboxamide-containing thiourea derivatives containing a phenylenediamine group, useful as inhibitors of herpes viruses

L31 ANSWER 9 OF 12 CAPLUS COPYRIGHT 2002 ACS
 TI Benzamide-containing aryl thiourea derivatives useful as inhibitors of herpes viruses

L31 ANSWER 10 OF 12 CAPLUS COPYRIGHT 2002 ACS
 TI Acetamide and substituted acetamide-containing aryl thiourea derivatives useful as inhibitors of herpes viruses

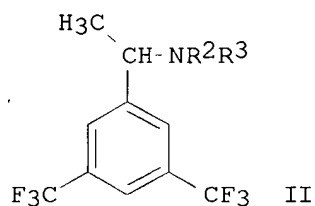
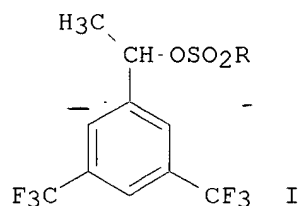
L31 ANSWER 11 OF 12 CAPLUS COPYRIGHT 2002 ACS
 TI Preparation of tryptophan derivatives as tachykinin antagonists

L31 ANSWER 12 OF 12 CAPLUS COPYRIGHT 2002 ACS
 TI Optically active amines. 34. Application of the benzene chirality rule to ring-substituted phenylcarbinamines and carbinols

=> d l31 4 ti fbib abs

L31 ANSWER 4 OF 12 CAPLUS COPYRIGHT 2002 ACS
 TI Preparation of optically active .alpha.-methyl-bis-3,5-(trifluoromethyl)benzylamine
 AN 2001:767496 CAPLUS
 DN 135:318319
 TI Preparation of optically active .alpha.-methyl-bis-3,5-(trifluoromethyl)benzylamine
 IN Ishii, Akio; Kuriyama, Masaru; Kanai, Masatomi; Hayami, Takashi
 PA Central Glass Co., Ltd., Japan
 SO Jpn. Kokai Tokkyo Koho, 11 pp.
 CODEN: JKXXAF
 DT Patent
 LA Japanese
 FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 2001294568	A2	20011023	JP 2000-112644	20000413
OS	CASREACT 135:318319; MARPAT 135:318319				
GI					



AB Title compd. is prepd. by sulfonylation of .alpha.-methyl-bis-3,5-(trifluoromethyl)benzyl alc. with RSO₂X (R = C1-6 alkyl, CmYnH2m+1-n, aryl; m = 0-8; n = 1-17; Y = F, Cl; X = F, Cl, RSO₂O), substitution of benzyl sulfonates I (R = same as above) with R¹NR²R³ (R¹ = H, alkali metal, alk. earth metal; R², R³ = H, arylalkyl, aryl, arylalkoxycarbonyl, OH, etc.), and deprotection of benzylamines II (R², R³ = same as above). (S)-.alpha.-methyl-bis-3,5-(trifluoromethyl)benzyl alc. was sulfonated with MeSO₂Cl in the presence of Et₃N in PhMe at 0.degree. for 1 h and reacted with benzylamine in DMF-PhMe at 65.degree. for 12 h to give 82% (R)-N-benzyl-.alpha.-methyl-bis-3,5-(trifluoromethyl)benzylamine, which was hydrogenated with H in the presence of P/C in EtOH at 56.degree. for 5 h to give 80% (R)-.alpha.-methyl-bis-3,5-(trifluoromethyl)benzylamine with 94.4% e.e.

=>

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COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
20.26	96.26

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
-1.24	-5.58

CA SUBSCRIBER PRICE

SESSION WILL BE HELD FOR 60 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 08:47:17 ON 25 APR 2002